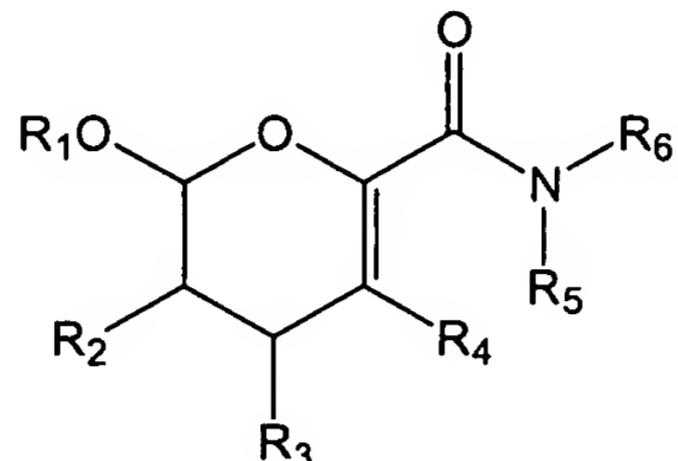


## AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions, and listings, of claims in the application:

1. **(Original)** A compound having the structure:



(I)

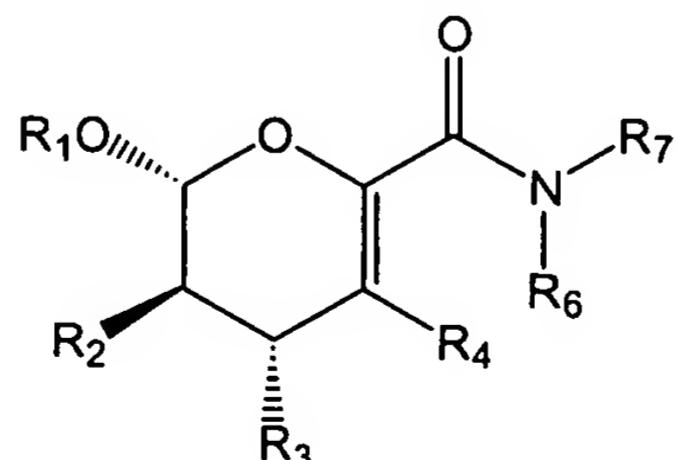
wherein R<sub>1</sub>-R<sub>4</sub> are each independently hydrogen or an aliphatic, heteroaliphatic, aryl, heteroaryl, alkylaryl or alkylheteroaryl moiety;

R<sub>5</sub> and R<sub>6</sub> are each independently hydrogen or an aliphatic, heteroaliphatic, aryl, heteroaryl, alkylaryl, or alkylheteroaryl moiety, and wherein R<sub>6</sub> and R<sub>7</sub>, taken together, may form a cyclic aliphatic, heteroaliphatic, aliphatic(aryl), heteroaliphatic(aryl), aliphatic(heteroaryl) or heteroaliphatic(heteroaryl) moiety, or an aryl or heteroaryl moiety;

wherein each of the foregoing aliphatic and heteroaliphatic moieties may be substituted or unsubstituted, cyclic or acyclic, saturated or unsaturated or linear or branched; and each of the foregoing aryl, heteroaryl, alkylaryl or alkylheteroaryl moieties may be substituted or unsubstituted; and

pharmaceutically acceptable derivatives thereof.

2. **(Original)** The compound of claim 1, wherein the compound has the structure (II):



(II)

wherein R<sub>1</sub>-R<sub>4</sub> are each independently hydrogen or an aliphatic, heteroaliphatic, aryl,

heteroaryl, alkylaryl or alkylheteroaryl moiety;

$R_5$  and  $R_6$  are each independently hydrogen or an aliphatic, heteroaliphatic, aryl, heteroaryl, alkylaryl, or alkylheteroaryl moiety, and wherein  $R_6$  and  $R_7$ , taken together, may form a cyclic aliphatic, heteroaliphatic, aliphatic(aryl), heteroaliphatic(aryl), aliphatic(heteroaryl) or heteroaliphatic(heteroaryl) moiety, or an aryl or heteroaryl moiety;

wherein each of the foregoing aliphatic and heteroaliphatic moieties may be substituted or unsubstituted, cyclic or acyclic, saturated or unsaturated or linear or branched; and each of the foregoing aryl, heteroaryl, alkylaryl or alkylheteroaryl moieties may be substituted or unsubstituted; and

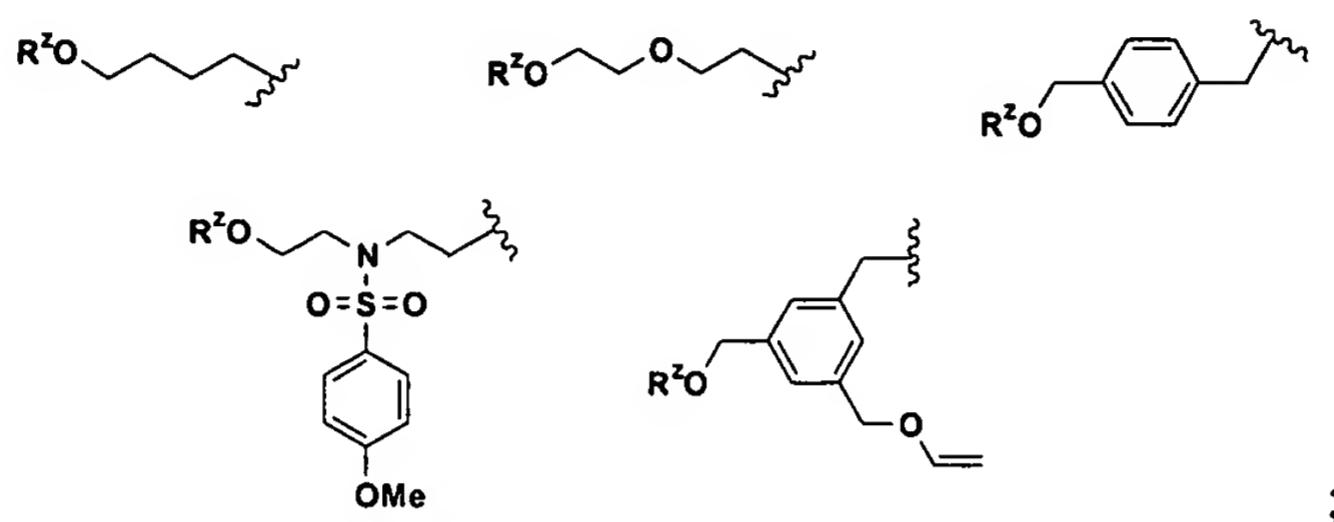
pharmaceutically acceptable derivatives thereof.

3. **(Original)** The compound of claim 1, wherein  $R^1$  is hydrogen or an alkyl, heteroalkyl, aryl or heteroaryl moiety substituted with  $Z$ , wherein  $Z$  is hydrogen,  $-(CH_2)_qOR^Z$ ,  $-(CH_2)_qSR^Z$ ,  $-(CH_2)_qN(R^Z)_2$ ,  $-(C=O)R^Z$ ,  $-(C=O)N(R^Z)_2$ , or an aliphatic, heteroaliphatic, aryl, heteroaryl, -(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety, wherein  $q$  is 0-4, and wherein each occurrence of  $R^Z$  is independently hydrogen, a protecting group, a solid support unit, or an aliphatic, heteroaliphatic, aryl, heteroaryl, -(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.

4. **(Original)** The compound of claim 3, wherein  $R^1$  is hydrogen, lower alkyl, a substituted or unsubstituted phenyl or -(lower alkyl)phenyl moiety,  $-(CH_2)_nOR^Z$ ,  $-[(CH_2)_nO]_mR^Z$ ,  $-(CH_2)_n-Ar-(CH_2)_mOR^Z$ ; wherein  $n$  and  $m$  are each independently integers from 1-6, Ar represents a substituted or unsubstituted aryl or heteroaryl moiety, and  $R^Z$  is independently hydrogen, a protecting group, a solid support unit, or an aliphatic, heteroaliphatic, aryl, heteroaryl, -(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of

the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.

5. **(Original)** The compound of claim 4, wherein  $R^1$  is hydrogen, ethyl, or has one of the structures:



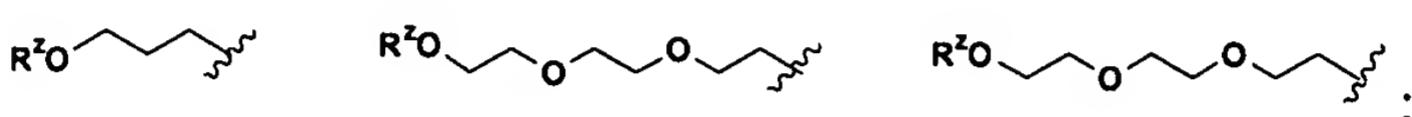
wherein  $R^z$  is as defined in claim 4.

6. **(Original)** The compound of claim 1, wherein  $R^2$  is hydrogen or an alkyl, heteroalkyl, aryl or heteroaryl moiety substituted with Z, wherein Z is hydrogen,  $-(CH_2)_qOR^z$ ,  $-(CH_2)_qSR^z$ ,  $-(CH_2)_qN(R^z)_2$ ,  $-(C=O)R^z$ ,  $-(C=O)N(R^z)_2$ , or an aliphatic, heteroaliphatic, aryl, heteroaryl, -(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety, wherein q is 0-4, and wherein each occurrence of  $R^z$  is independently hydrogen, a protecting group, a solid support unit, or an aliphatic, heteroaliphatic, aryl, heteroaryl, -(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.

7. **(Original)** The compound of claim 6, wherein  $R^2$  is hydrogen, lower alkyl, a substituted or unsubstituted phenyl or -(lower alkyl)phenyl moiety,  $-(CH_2)_nOR^z$ ,  $-[(CH_2)_nO]_mR^z$ ,  $-(CH_2)_n-Ar-(CH_2)_mOR^z$ ; wherein n and m are each independently integers from 1-6, Ar represents a substituted or unsubstituted aryl or heteroaryl moiety, and  $R^z$  is independently hydrogen, a protecting group, a solid support unit, or an aliphatic, heteroaliphatic, aryl, heteroaryl, -

(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.

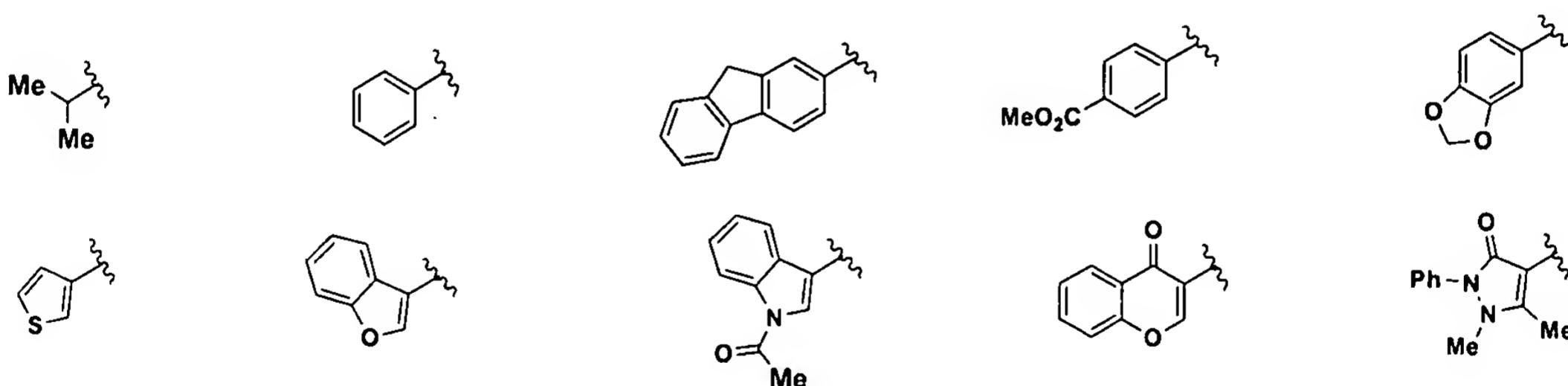
8. (Original) The compound of claim 6, wherein  $R^2$  is hydrogen or has one of the structures:



wherein  $R^z$  is as defined in claim 6.

9. (Original) The compound of claim 1, wherein  $R^3$  is an alkyl, heteroalkyl, aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.

10. (Original) The compound of claim 9, wherein  $R^3$  has one of the structures:



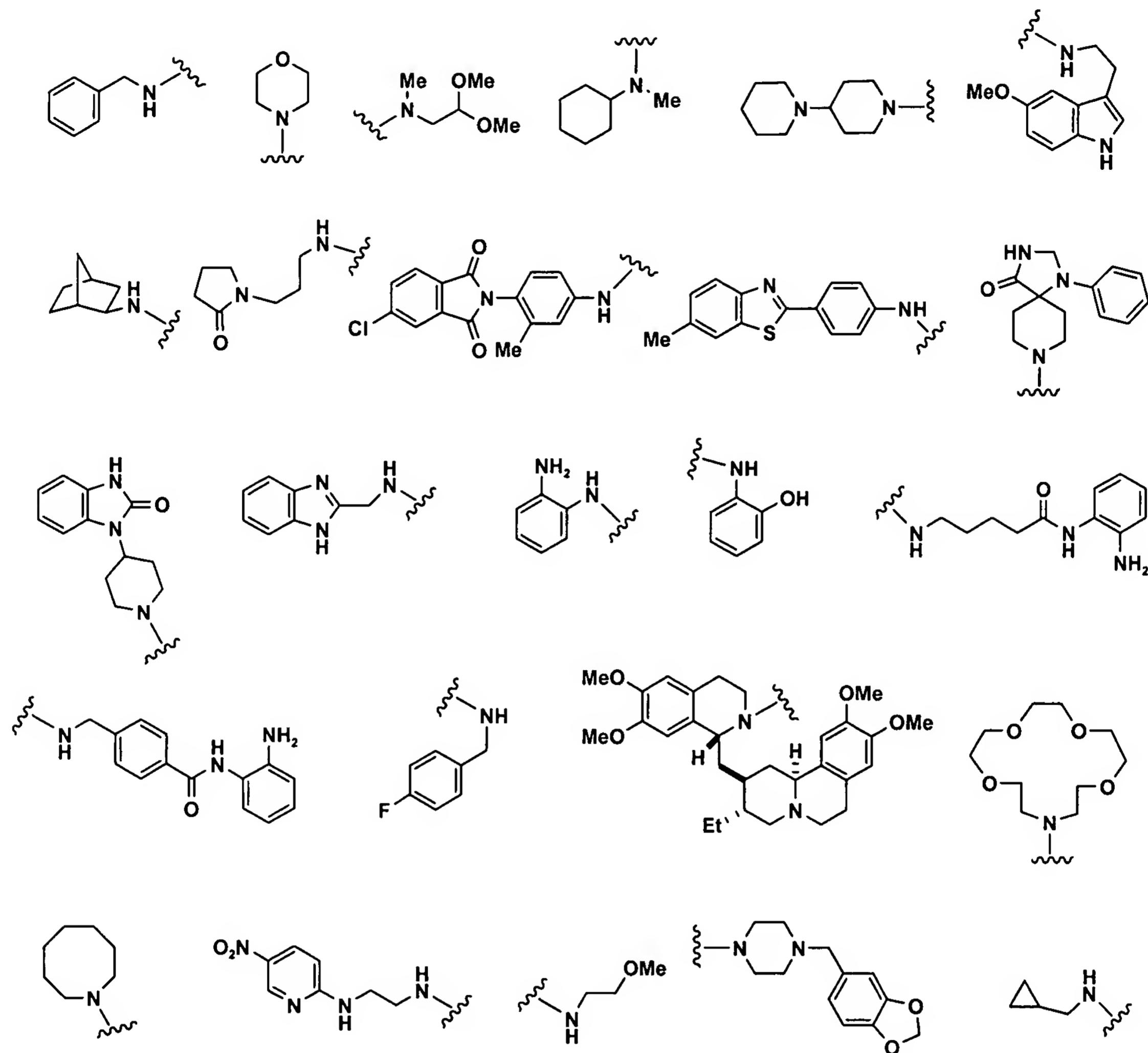
11. (Original) The compound of claim 1, wherein R<sup>4</sup> is hydrogen or an alkyl, heteroalkyl, aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of

the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or –(heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.

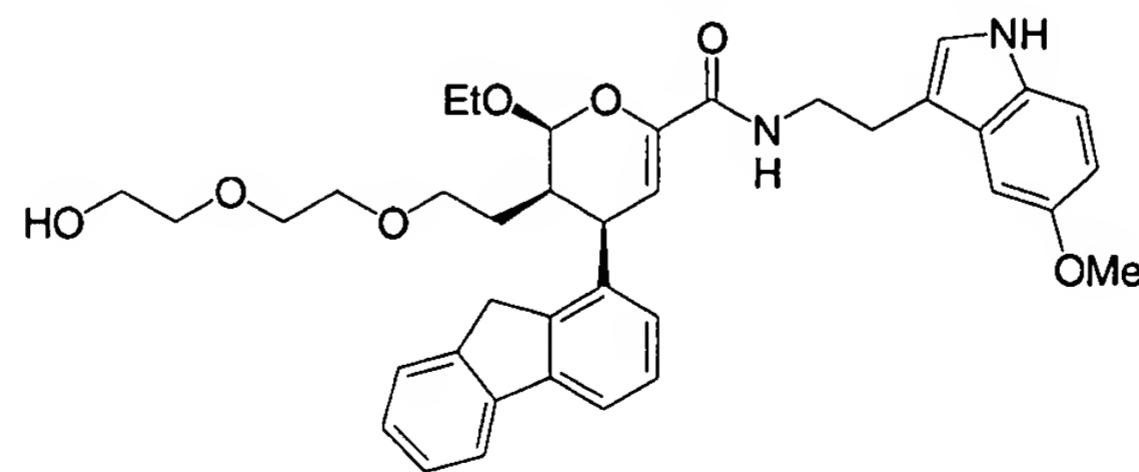
12. **(Original)** The compound of claim 11, wherein R<sup>4</sup> is hydrogen alkyl or heteroalkyl.

13. **(Original)** The compound of claim 1, wherein R<sup>5</sup> and R<sup>6</sup> are each independently hydrogen or an alkyl, heteroalkyl, aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or –(heteroalkyl)heteroaryl moiety; or wherein R<sup>5</sup> and R<sup>6</sup>, taken together, form a substituted or unsubstituted, saturated or unsaturated cyclic moiety comprising 5-12 carbon atoms, 0-5 oxygen atoms, 0-5 sulfur atoms and 1-5 nitrogen atoms; and wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or –(heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.

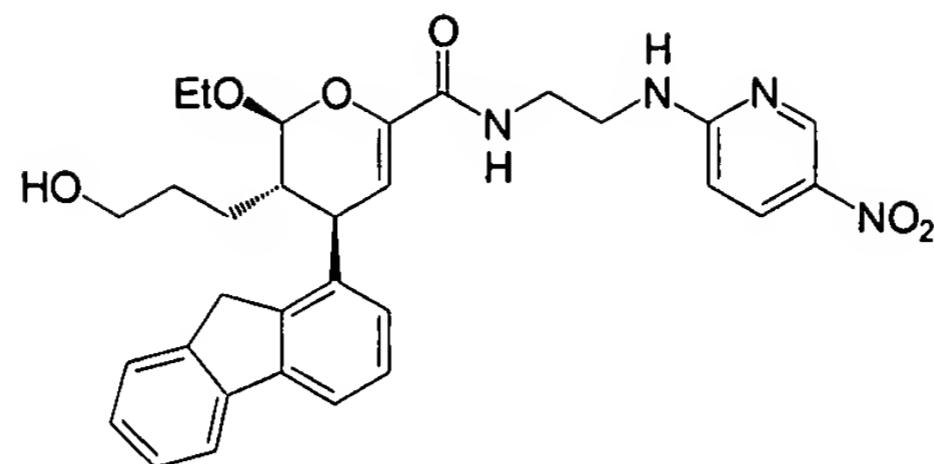
14. **(Original)** The compound of claim 1, wherein –NR<sup>5</sup>R<sup>6</sup> is one of the following the structures:



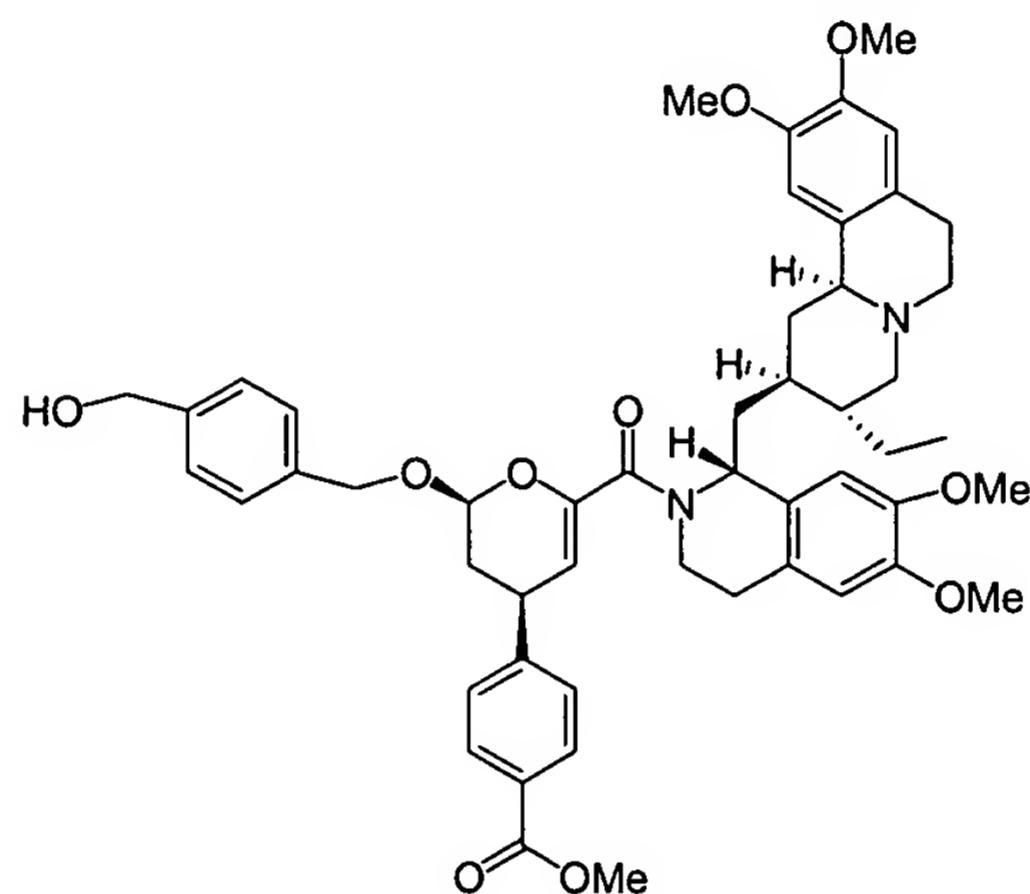
15. (Original) The compound of claim 1 having the structure:



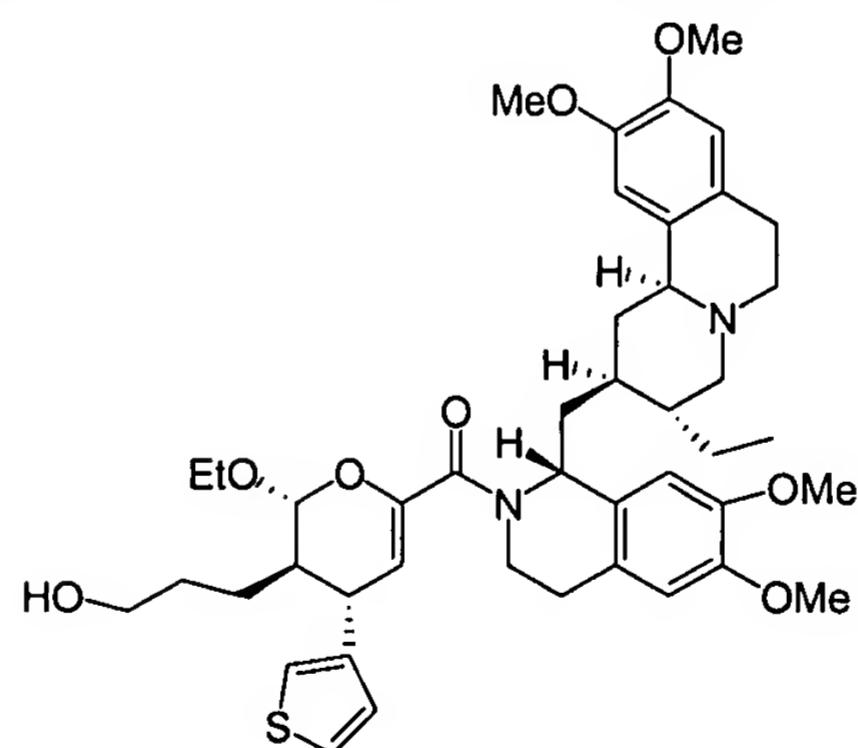
16. (Original) The compound of claim 1 having the structure:



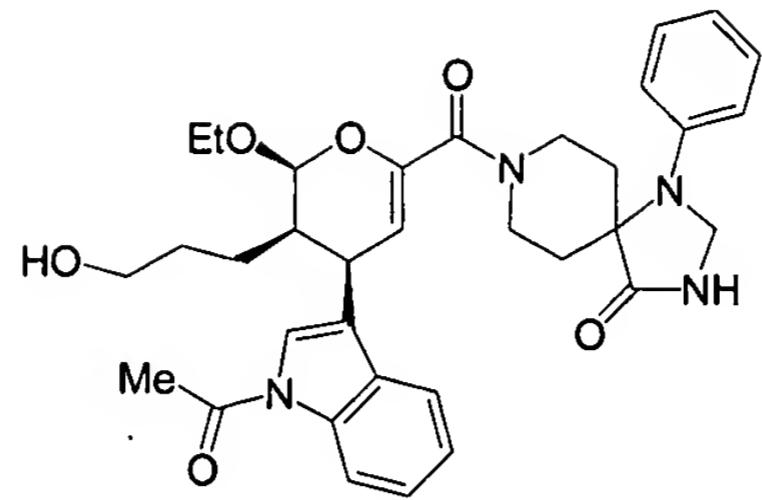
17. **(Original)** The compound of claim 1 having the structure:



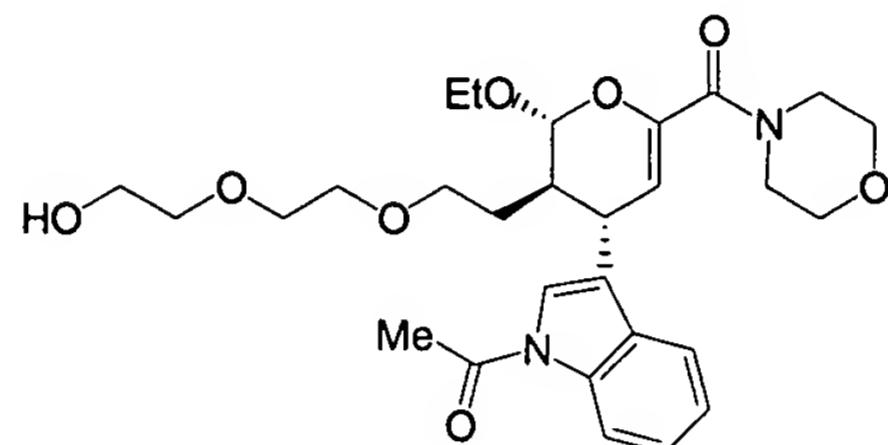
18. **(Original)** The compound of claim 1 having the structure:



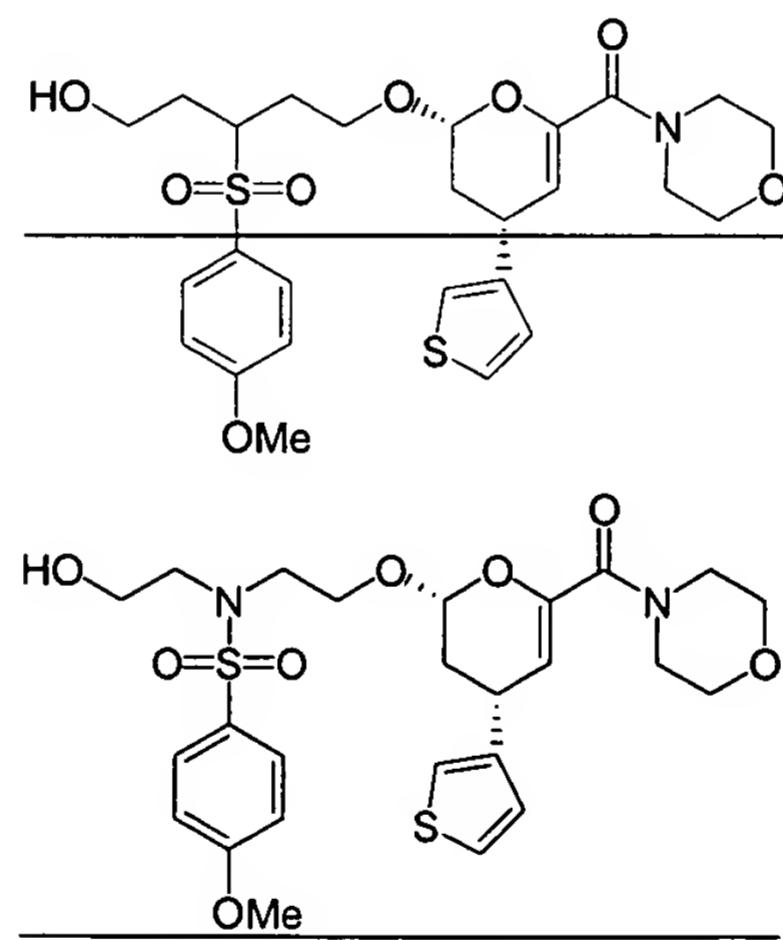
19. **(Original)** The compound of claim 1 having the structure:



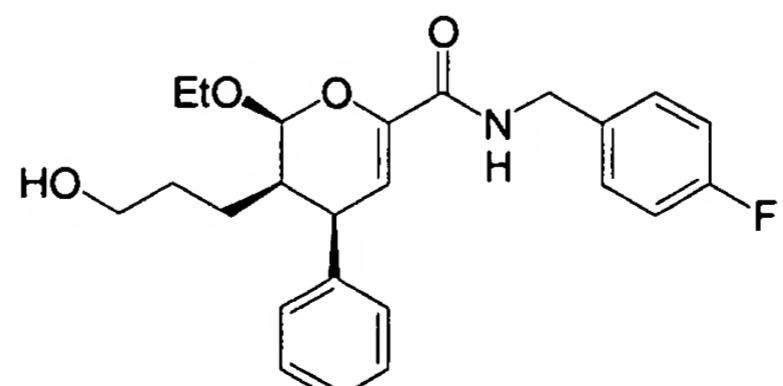
20. **(Original)** The compound of claim 1 having the structure:



21. **(Currently Amended)** The compound of claim 1 having the structure:



22. **(Original)** The compound of claim 1 having the structure:



23. **(Original)** A collection of compounds comprising two or more of the compounds of claim 1 or 2.
24. **(Original)** The collection of claim 23, wherein the collection is provided in array format.
25. **(Original)** The collection of claim 23, wherein the collection is provided in array format on a glass slide.
26. **(Original)** The collection of claim 23, wherein the collection comprises at least 100 compounds.
27. **(Original)** The collection of claim 23, wherein the collection comprises at least 1,000 compounds.
28. **(Original)** The collection of claim 23, wherein the collection comprises at least 2,000 compounds.
29. **(Original)** The collection of claim 23, wherein the collection comprises at least 10, 000 compounds.
30. **(Original)** A pharmaceutical composition comprising:  
a compound of any one of claims 1, 2, 5, 8, 10, 14, and 15-22; and  
a pharmaceutically acceptable carrier.

**Claims 31-34 Canceled**

35. **(Original)** A method for inhibiting a kinesin activity comprising contacting a cell with a compound of any one of claims 1, 2, 5, 8, 10, 14, and 15-22.
36. **(Original)** The method of claim 35, wherein the kinesin is Eg5.

37. **(Original)** A method for treating a proliferative disorder comprising: administering to a subject in need thereof a therapeutically effective amount of a compound of any one of claims 1, 2, 5, 8, 10, 14, and 15-22.
38. **(Original)** The method of claim 37, wherein the proliferative disorder is cancer.
39. **(Original)** The method of claim 37, further comprising administering an additional therapeutic agent.